

Text Classification with Multi-Layer Perceptrons

2024–25

Ion Androutsopoulos http://www.aueb.gr/users/ion/ These slides are partly based on material from the books:

- Artificial Intelligence A Modern Approach by S. Russel and P. Norvig, 2nd edition, Prentice Hall, 2003,
- Artificial Intelligence by I. Vlahavas et al., 3rd edition, University of Macedonia Press, 2006 (in Greek).
- *Machine Learning* by T. Mitchell, McGraw-Hill, 1997.

Contents

- Natural and artificial neural networks (NNs).
- Perceptrons, training them with SGD, limitations.
- Multi-Layer Perceptrons (MLPs) and backpropagation.
- MLPs for text classification, regression, window-based token classiffication (e.g., for POS tagging, NER).
- Dropout, batch and layer normalization.
- Pre-training word embeddings with Word2Vec.
- Advice for training large neural networks.

Natural neural networks



- Neuron: cell of the brain.
 - Cell body or soma: the main part, includes the nucleus.
 - Dendrites receive signals from other neurons.
 - Axon: transmits a single output to other neurons. Often much longer than the diameter of the soma.
 - Synapses: axon-dendrite interfaces, whose conductivities vary.
- Neural network: network of many neurons.

Artificial neural networks

- Artificial neuron:
 - Input: real variables.
 - Input weights: real variables (roughly synapses).
 - Soma: computes the weighted sum of the inputs, then applies an activation function to the sum.



Activation functions





ReLU and variants are popular choices.

and equations from Left image from http://datascience.stackexchange.com/questions/5706/. Right image https://fleuret.org/francois/ The Deep 1 Learning, by F. Fleuret

Perceptron

- A single neuron, originally with sign activation function.
 - Equivalently, step activation function.
 - The Perceptron can be generalized (done later), to use a sigmoid or other activation function.
 - We can use several Perceptrons (e.g., to recognize a letter each).



We usually think of it as an algorithm, but it was a machine originally! The perceptron is a machine



Slide from the presentation "Multilayer Neural Networks" of L. Bottou at the Deep Learning Summer School 2015 (http://videolectures.net/deeplearning2015_montreal/).

Perceptrons as logical gates

true: 1, false: -1



Two-level XOR implementation





- Like logistic regression, a single **Perceptron** can learn only **linear separators** (exercise). For **non-linear separators**, we need **Multi-Layer Perceptrons**.
- An MLP with 1 hidden layer can compute (in principle, also learn) any mapping between discrete spaces with finite dimensions.
- For continuous spaces, very roughly speaking: with 1 hidden layer we can compute (in principle, also learn) almost any bounded continuous function; with 2 hidden layers we can compute almost any bounded function.
- But we may need a very large, unknown number of neurons in the hidden layers, we may end up memorizing the training dataset (exercise), and we may not actually manage to learn (find) the target function.

Perceptron's original learning algorithm

- 1. Start with random weights \vec{w} .
- 2. Set $i \leftarrow 1$ and $s \leftarrow 0$.
- 3. Let $t^{(i)}$ be the correct output for the *i*-th training instance and $o^{(i)}$ the current output for that instance.
- 4. Set $s \leftarrow s + E_i(\vec{w})$, with: $E_i(\vec{w}) = 1/2 \cdot [t^{(i)} o^{(i)}]^2$
- 5. Update the weights: $w_l \leftarrow w_l + \eta \cdot (t^{(i)} o^{(i)}) \cdot x_l^{(i)}$
- 6. If there is a next training instance, set $i \leftarrow i+1$ and go to step 3.
- 7. If *s* had not converged and max number of scans (epochs) of training data not exceeded go to step 2.
- In the simplest case, η is a small positive constant.

Perceptron with sigmoid or other Φ

 More generally, when using a Perceptron with activation function Φ instead of step/sign:

$$w_l \leftarrow w_l + \eta \left(\sum_l w_l x_l^{(i)} \right) \cdot \left(t^{(i)} - \Phi \left(\sum_l w_l x_l^{(i)} \right) \right) \cdot x_l^{(i)}$$

- For sigmoid $\Phi(S) = \frac{1}{1 + e^{-S}}, \Phi'(S) = \Phi(S) \cdot (1 \Phi(S)).$
- And since $o^{(i)} = \Phi\left(\sum_{l} w_{l} x_{l}^{(i)}\right)$, the weights update rule becomes:

$$w_l \leftarrow w_l + \eta \cdot o^{(i)} \cdot (1 - o^{(i)}) \cdot (t^{(i)} - o^{(i)}) \cdot x_l^{(i)}$$



Derivation of the update rule

$$E_{i}(\vec{w}) = \frac{1}{2} (t^{(i)} - o^{(i)})^{2} = \frac{1}{2} (t^{(i)} - \Phi\left(\sum_{l=1}^{n} w_{l} x_{l}^{(i)}\right))^{2}$$

$$\nabla E_{i}(\vec{w}) = \left\langle \frac{\partial E_{i}(\vec{w})}{\partial w_{0}}, \dots, \frac{\partial E_{i}(\vec{w})}{\partial w_{l}}, \dots, \frac{\partial E_{i}(\vec{w})}{\partial w_{n}} \right\rangle$$

$$\frac{\partial E_{i}}{\partial w_{l}} = (t^{(i)} - \Phi\left(\sum_{l} w_{l} x_{l}^{(i)}\right)) \cdot \frac{\partial (t^{(i)} - \Phi\left(\sum_{l} w_{l} x_{l}^{(i)}\right))}{\partial w_{l}} = -(t^{(i)} - o^{(i)}) \cdot \Phi'(\sum_{l} w_{l} x_{l}^{(i)}) \cdot x_{l}^{(i)}$$
Hence: $\nabla E_{i}(\vec{w}) = -(t^{(i)} - o^{(i)}) \cdot \Phi'(\sum_{l} w_{l} x_{l}^{(i)}) \cdot \left\langle x_{1}^{(i)}, \dots, x_{n}^{(i)} \right\rangle$

$$= -(t^{(i)} - o^{(i)}) \cdot \Phi'(\vec{w} \cdot \vec{x}^{(i)}) \cdot \vec{x}^{(i)}$$

Weights update rule:

$$\vec{w} \leftarrow \vec{w} - \eta \cdot \nabla E_i(\vec{w}) = \vec{w} + \eta \cdot (t^{(i)} - o^{(i)}) \cdot \Phi'(\vec{w} \cdot \vec{x}^{(i)}) \cdot \vec{x}^{(i)}$$

For each weight: $w_l \leftarrow w_l + \eta \cdot (t^{(i)} - o^{(i)}) \cdot \Phi'(\vec{w} \cdot \vec{x}^{(i)}) \cdot x_l^{(i)}$



Backpropagation Original image from the book of Vlahavas et al. hidden layer 4 **W**₁₄ **W**48 **W**24 **W**34 **X**1 **W**15 **W**49 **y**8 8 compute the output for each input **X**2 **W**35 2 W68 backpropagate derivatives to adjust weights **y**9 9 **W**36 **W**78 **W**₁₇ **X**3 3 W27 **W**79 **W**37 7 output layer input layer

Backpropagation

- Initialize all the weights to small random values.
 - \circ E.g., sample from a **zero-centered Gaussian** with small σ .
 - Better initializations exist (see DL course).
 - Normalize the features too (see "Important tricks" of Part 2).
- In each epoch, for each training example (or mini-batch):
 - Compute the **output** $< o_1, o_2, \ldots >$ for the training example.
 - For each weight w_{ij} , compute $\frac{\partial E}{\partial w_{ij}}$, where *E* the loss on the training example. We compute derivatives right to left.
 - Update each weight as: $w_{ij} \leftarrow w_{ij} \eta \cdot \frac{\partial E}{\partial w_{ij}}$, i.e., for all the weights together: $W \leftarrow W \eta \cdot \nabla_W E$
 - Hence, we use **SGD** (or variants). **No guarantee** SGD will find the **best solution**, but it (often) works in practice!

Example of computation graph



Example and figure from Stanford's "CNNs for Visual Recognition" (2016, F.-F. Li, A. Karpathy, J. Johnson) http://cs231n.github.io/optimization-2/

- Forward pass: $\langle x, y, z \rangle = \langle -2, 5, -4 \rangle, q = 3, f = -12$
- Imagine we wish to **minimize** *f* using **SGD**.
 - In a more realistic scenario, f would be a loss function, and $\langle x, y, z \rangle$ the weights vector.

 $\begin{bmatrix} x \\ y \\ z \end{bmatrix} \leftarrow \begin{bmatrix} x \\ y \\ z \end{bmatrix} - \eta \nabla f(x, y, z) = \begin{bmatrix} x \\ y \\ z \end{bmatrix} - \eta \begin{vmatrix} \overline{\partial x} \\ \overline{\partial f} \\ \overline{\partial y} \\ \overline{\partial f} \end{vmatrix} \xrightarrow{\frac{\partial f}{\partial x}, \frac{\partial f}{\partial y}, \frac{\partial f}{\partial z}}$

We need:

Backpropagation in the graph

-12



Example and figure from Stanford's "CNNs for Visual Recognition" (2016, F.-F. Li, A. Karpathy, J. Johnson) http://cs231n.github.io/optimization-2/

- Backpropagation: We compute derivatives right to left. $\circ \frac{\partial f}{\partial f} = 1$ by definition.
 - $\circ \frac{\partial f}{\partial q} = \dot{z}$. And for this $\langle x, y, z \rangle$ input, z = -4.
 - During the forward pass, we need to save the outputs of all the nodes (e.g., here we need the value of z).

Backpropagation in the graph



-4 3 5 y -12 *

Example and figure from Stanford's "CNNs for Visual Recognition" (2016, F.-F. Li, A. Karpathy, J. Johnson) http://cs231n.github.io/optimization-2/

Backpropagation: We compute derivatives right to left. ·a f (

$$\frac{\partial f}{\partial t} = 1$$
 by definition.

$$\circ \frac{\partial f}{\partial q} = z$$
. And for this $\langle x, y, z \rangle$ input, $z = -4$.

$$\circ \frac{\partial f}{\partial z} = q$$
. And for this $\langle x, y, z \rangle$ input, $q = 3$.

$$\circ \frac{\partial f}{\partial y} = \frac{\partial f}{\partial q} \frac{\partial q}{\partial y} = \frac{\partial f}{\partial q} \cdot 1. \text{ And here } \frac{\partial f}{\partial q} \text{ is } -4.$$
$$\circ \frac{\partial f}{\partial x} = \frac{\partial f}{\partial q} \frac{\partial q}{\partial x} = \frac{\partial f}{\partial q} \cdot 1. \text{ And here } \frac{\partial f}{\partial q} \text{ is } -4.$$

incoming gradient local gradient



class PlusGate: forward(x, y): return x+y backward($\frac{\partial f}{\partial a}$): return $< \frac{\partial f}{\partial q}, \frac{\partial f}{\partial q} >$



class StarGate: forward(q, z): return q * z backward($\frac{\partial f}{\partial w}$): return $< \frac{\partial f}{\partial w} \cdot z, \frac{\partial f}{\partial w} \cdot q >$

More compact notation of NNs



More compact notation of NNs

$$\begin{array}{c} x_{1} & w_{1,1} & n_{1} \\ w_{2,1} & w_{m,1} \\ x_{2} & w_{m,1} \\ x_{3} & w_{m,k} \\ & & & & \\ & & & \\ &$$

More compact notation of NNs



$$\vec{o}^{(1)} = \begin{bmatrix} o_{1,1} \\ o_{1,2} \\ \dots \\ o_{1,k_1} \end{bmatrix} = \Phi(\vec{s}^{(1)}) = \Phi(W^{(1)}\vec{x})$$







Classification example



Softmax



$$= \begin{bmatrix} \exp(s_{2,1}) \\ \overline{\sum_{j=1}^{k_2} \exp(s_{2,j})} \\ \exp(s_{2,2}) \\ \overline{\sum_{j=1}^{k_2} \exp(s_{2,j})} \\ \dots \\ \exp(s_{2,k_2}) \\ \overline{\sum_{j=1}^{k_2} \exp(s_{2,j})} \end{bmatrix}$$

Softmax also **moves the largest of its inputs towards 1** and the other inputs towards 0. Intuitively a **soft argmax**!

Classification example



Classification example – more compact



Or as a **computation graph**:



Extracting contract elements

THIS AGREEMENT is made the 15th day of October 2009 (The "Effective Date") BETWEEN:

Sugar 13 Inc., a corporation whose office is at James House,
 42-50 Bond Street, London, EW2H TL ("Sugar");

(2) **E2 UK Limited**, a limited company whose registered office is at 260 Bathurst Road, Yorkshire, SL3 4SA (**"Provider"**).

RECITALS:

A. The Parties wish to enter into a framework agreement which will enable Sugar, from time to time, to [...]B. [...]

NO THEREFORE IT IS AGREED AS FOLLOWS:

ARTICLE I - DEFINITIONS

"Sugar" shall mean: Sugar 13 Inc.

"Provider" shall mean:	E2 UK Limited
------------------------	---------------

"1933 Act" shall mean: Securities Act of 1933

ARTICLE II - TERMINATION

The Service Period will be for five (5) years from the Effective Date (The "Initial Term"). The agreement is considered to be terminated in October 16, 2014.

ARTICLE III - PAYMENT - FEES

During the service period monthly payments should occur. The estimated fees for the Initial Term are £154,800.

ARTICLE IV - GOVERNING LAW

This agreement shall be governed and construed in accordance with the Laws of England & Wales. Each party hereby irrevocably submits to the exclusive jurisdiction of the courts sitting in Northern London.

IN WITNESS WHEREOF, the parties have caused their respective duly authorized officers to execute this Agreement.

BY: George Fake Authorized Officer Sugar 13 Inc.

BY: Olivier Giroux CEO E2 UK LIMITED

Identify start/end dates, duration, contractors, amount, legislations refs, jurisdiction etc. Similar to Named Entity Recognition (NER).

I. Chalkidis, I. Androutsopoulos, A. Michos, "Extracting Contract Elements", ICAIL 2017, http://nlp.cs.aueb.gr/pubs/icail2017.pdf.

Window-based NER example *i*-th word of the text being classified 3-word **window** (often larger) yesterday language tech announced that... $\vec{x}_{i-1} = \begin{bmatrix} 1\\0\\0\\0\\...\\0 \end{bmatrix} \quad \vec{x}_i = \begin{bmatrix} 0\\0\\1\\0\\...\\0 \end{bmatrix} \quad \vec{x}_{i+1} = \begin{bmatrix} 0\\0\\0\\0\\...\\0 \end{bmatrix} \quad \textbf{i-1} \quad \textbf{i-hot vectors } (|V| \times 1) \text{ of the window. } (|V| \text{ is the vocabulary size}).$ words in the window. (d $\vec{e}_{i-1} = \begin{bmatrix} 1.8\\ 2.3\\ -1.4\\ 3.7\\ ...\\ -1.1 \end{bmatrix} \vec{e}_i = \begin{bmatrix} 2.4\\ -3\\ 9.3\\ 5.1\\ ...\\ 3.9 \end{bmatrix} \vec{e}_{i+1} = \begin{bmatrix} 2.2\\ 3.8\\ 1.2\\ -6.4\\ ...\\ 7.1 \end{bmatrix}$ Is the **dimensionality** of the word embeddings). Is the **dimensionality** of the word embeddings. Let \boldsymbol{E} be a matrix $(d \times |V|)$ that **contains** all the **embeddings** of the **vocabulary** as **columns**. Then: $\vec{e}_{i-1} = F\vec{v} \cdot \vec{c} - F\vec{v}$ is the dimensionality of

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Window-based NER example



We learn $W^{(1)}$, $W^{(2)}$ with **backpropagation**. We can also learn (or modify) the **word embeddings** *E* during **backpropagation**! But when we don't have large training datasets (e.g., corpus manually annotated with B-I-O tags), it may be better to use **pre-trained embeddings**, which can be obtained from large non-annotated corpora (e.g., via Word2Vec, GloVe, to be discussed).

We can use the same window-based approach for **POS-tagging**, **chunking**, ...

Cross-entropy loss





Dropout

- For each training instance (or mini-batch), we drop (remove) each neuron of the layer where dropout is applied with probability $p_{drop} = 1 - p_{keep}$.
 - Helps the neural net **avoid relying too much on particular neurons** (or inputs). A form of **regularization**. Works well!
 - Gradients also do not flow through dropped neurons.
 - Alternative explanation: we train an **ensemble** of networks, containing **all the pruned networks** that dropout creates.
- **During testing**, we **multiply the output** of each neuron (of the layer where dropout was applied) by p_{keep} , so that the neuron's **expected output value** will be **as in training**.
 - Or we divide the output by p_{keep} during training instead.
 - We **don't drop neurons during testing** (only during training).

Batch normalization



At each layer, instead of: $s_j = \sum_{i=1}^{m} w_{i,j} x_i$

we use:

$$\bar{s}_j = \frac{g_j}{\sigma_j} (s_j - \mu_j) + b_j$$

- μ_j , σ_j are the **mean** and **std. dev. of** s_j in the **mini-batch**.
- g_j, b_j are **learned** parameters (constant after training).

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• Φ now applied to \bar{s}_j .

See <u>https://arxiv.org/pdf/1607.06450.pdf</u> for **batch vs. layer normalization**. The latter is better for RNNs (next part), where layers are time-steps.

Layer normalization



At each layer, instead of: $s_j = \sum_{i=1}^{m} w_{i,j} x_i$

we use:

$$\bar{s}_j = \frac{g_j}{\sigma} (s_j - \mu) + b_j$$

- μ , σ are the **mean** and **std. dev. of s₁**, ..., **s_k** in the <u>layer</u>.
- g_j , b_j are **learned** parameters (constant after training).

• Φ applied to \bar{s}_j .

With dropout, batch/layer normalization, residuals (to be discussed) and other additions, strictly speaking we no longer have an "MLP". Some people prefer "Feed Forward Neural Network" (FFNN), but "MLP" still often used as synonym.

Pretraining word embeddings with Word2Vec (skip-gram version)

- Every word w of the vocabulary has two vectors: $\vec{w}^{(in)}, \vec{w}^{(out)}$
 - The vectors are randomly initialized. We learn them.
 - For every token w_t at position t of a corpus and every position t + i ($i \neq 0$) within a window [t c, t + c] around position t:

$$w_{t+i} = \cdots?$$

$$w_{t+i} = "starring"?$$

$$w_{t+i} = "directed"?$$

$$w_t = "film"$$

$$w_{t+i} = "famous"?$$

$$w_{t+i} = t + i$$

• We want to be able to predict which vocabulary word occurs at position t + i by multiplying $\vec{w}_t^{(in)}$ and $\vec{w}_{t+i}^{(out)}$.

Word2Vec (skip-gram version)



Word2Vec (skip-gram version)

• We learn the $\vec{w}^{(in)}$, $\vec{w}^{(out)}$ by maximizing the probability assigned to the w_{t+i} that actually occurs at each position t + i, i.e., we maximize the likelihood of the correct predictions:

$$\left\langle \hat{E}^{(in)}, \hat{E}^{(out)} \right\rangle = \operatorname*{argmax}_{\left\langle E^{(in)}, E^{(out)} \right\rangle} \sum_{t=1+c}^{T-c} \sum_{-c \le i \le c, i \ne 0} \ln P(w_{t+i}|w_t)$$

- Matrices $E^{(in)}$, $E^{(out)}$ contain in their columns all the *in* and *out* vectors (word embeddings) of all vocabulary words.
- *T* is the corpus size, *t* is the center of the sliding window.
- For each *t* value, we get 2*c* training examples.
- For batch gradient ascent, we would do steps: $\langle \hat{E}^{(in)}, \hat{E}^{(out)} \rangle \leftarrow \langle \hat{E}^{(in)}, \hat{E}^{(out)} \rangle + \eta \nabla L$
- In practice, we use SGD (or variants), i.e., we use the likelihood L_i of a mini-batch of training examples (e.g., all 2*c* of a window): $\langle \hat{E}^{(in)}, \hat{E}^{(out)} \rangle \leftarrow \langle \hat{E}^{(in)}, \hat{E}^{(out)} \rangle + \eta \nabla L_i$



Word2Vec (skip-gram with negative sampling)



Word2Vec (skip-gram with negative sampling)



Loss as a function of epochs



Figure from the **recommended book "Deep Learning with Python"** by F. Chollet, Manning Publications, 2nd edition. The 1st edition is freely available. <u>https://www.manning.com/books/deep-learning-with-python</u> <u>https://www.manning.com/books/deep-learning-with-python-second-edition</u>

Practical advice for training deep NNs

- Check simple baselines: (e.g., majority, random, ...)
 - If you can't beat them, you may have **bugs**, **data problems**, ...
 - Look at how data are **tokenized**, **pre-processed**, ...
 - Examine misclassification errors (e.g., extreme/frequent cases).
- Get the training and validation loss to start falling:
 Tune the learning rate and the mini-batch size.
 Use appropriate models (e.g., for sequences, images, ...).
- Reach the overfitting behavior of the previous slide.
 - The **training and validation loss** (or metric) **both fall up to a point**, then the **training loss continues to improve** ideally reaching **near zero**, the **validation loss deteriorates**.
 - Increase capacity (e.g., layers, neurons per layer), ...
- Then dropout, early stopping, batch/layer norm, ...
- Check Chollet's Chapter 5 for more advice...

Regularizing a high-capacity model

Figure 5.21 Effect of dropout on validation loss



Figure from the **recommended book "Deep Learning with Python"** by F. Chollet, Manning Publications, 2nd edition. The 1st edition is freely available. <u>https://www.manning.com/books/deep-learning-with-python</u> https://www.manning.com/books/deep-learning-with-python-second-edition

Recommended reading

- M. Surdeanu and M.A. Valenzuela-Escarcega, *Deep Learning for Natural Language Processing: A Gentle Introduction*, Cambridge Univ. Press, 2024.
 - Chapters 5–9.
 - <u>https://clulab.org/gentlenlp/text.html</u>
 - Also available at AUEB's library.
- Y. Goldberg, *Neural Network Models for Natural Language Processing*, Morgan & Claypool Publishers, 2017.
 Mostly chapters 3–5 and 10. Available at AUEB's library.
- Jurafsky & Martin's, *Speech and Language Processing* is being revised (3rd ed.) to include Deep Learning methods.
 - <u>http://web.stanford.edu/~jurafsky/slp3/</u>



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Other recommended resources

- For an introduction to Keras/Tensorflow and practical DL for NLP and vision, see F. Chollet's *Deep Learning in Python*, Manning Publications, 1st edition, 2017.
 - The 1st edition is freely available and sufficient for this course. <u>https://www.manning.com/books/deep-learning-with-python</u>
 - 2nd edition (2022) now available, requires payment. Highly recommended.
- Useful maths background: T. Parr και J. Howard, *The Matrix Calculus You Need for Deep Learning*.

o https://explained.ai/matrix-calculus

- PyTorch tutorials: <u>https://pytorch.org/tutorials/</u>
- C. Manning's (Stanford) *NLP with Deep Learning* course.
 <u>http://web.stanford.edu/class/cs224n/</u>. Videos on YouTube.
- See also the recommended books of Part 0 (Introduction).



